



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:38 PM UTC

PDB ID : 3CL9 / pdb_00003cl9
Title : Structure of bifunctional TcDHFR-TS in complex with MTX
Authors : Schormann, N.; Senkovich, O.; Chattopadhyay, D.
Deposited on : 2008-03-18
Resolution : 3.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

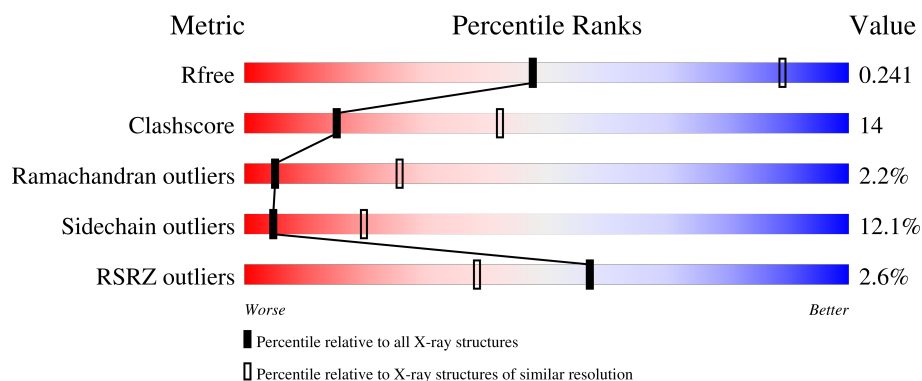
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	521	

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 4157 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

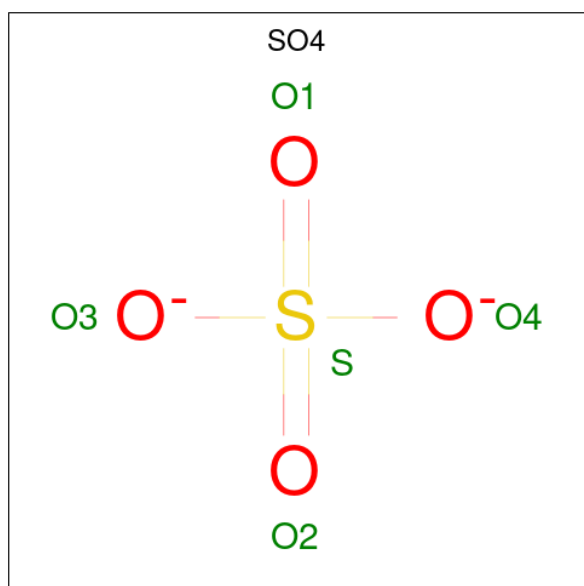
- Molecule 1 is a protein called Bifunctional dihydrofolate reductase-thymidylate synthase (DHFR-TS).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	505	4015	2549	709	739	18	0	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	32	ARG	HIS	SEE REMARK 999	UNP Q27793
A	55	VAL	LEU	SEE REMARK 999	UNP Q27793
A	137	GLN	ARG	SEE REMARK 999	UNP Q27793

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0

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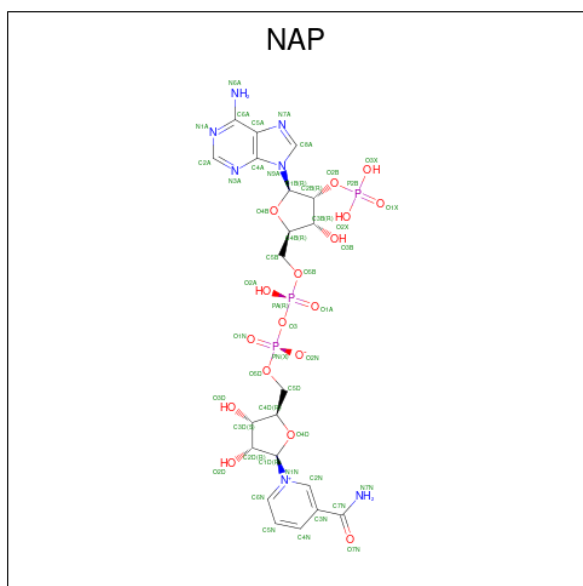
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

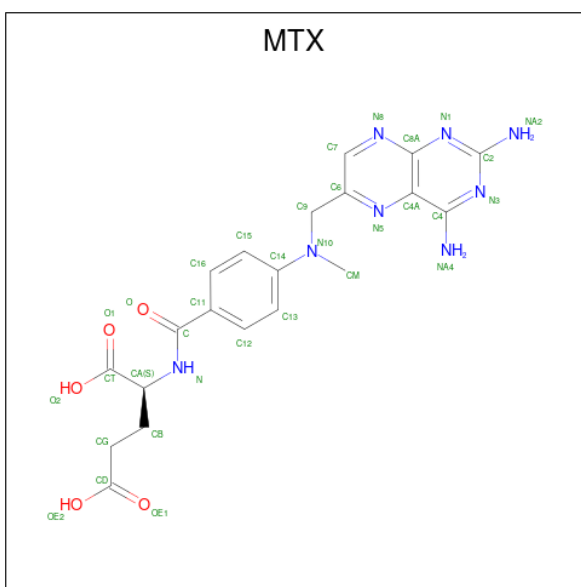
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	3	Total Cl 3 3	0	0

- Molecule 4 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: $\text{C}_{21}\text{H}_{28}\text{N}_7\text{O}_{17}\text{P}_3$).



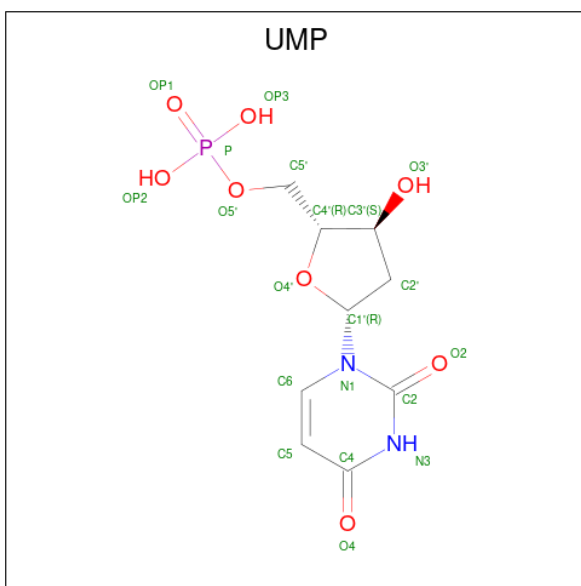
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 48	C 21	N 7	O 17	P 3	0	0

- Molecule 5 is METHOTREXATE (CCD ID: MTX) (formula: $C_{20}H_{22}N_8O_5$).



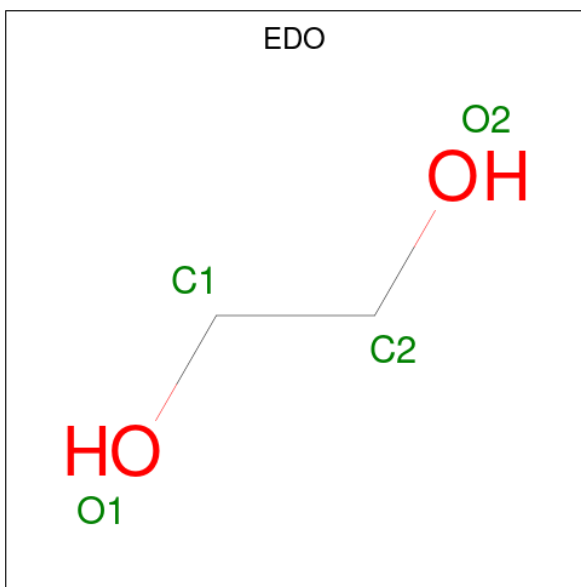
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			33	20	8	5		

- Molecule 6 is 2'-DEOXYURIDINE 5'-MONOPHOSPHATE (CCD ID: UMP) (formula: $C_9H_{13}N_2O_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			20	9	2	8	1		

- Molecule 7 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		
7	A	1	Total	C	O	0	0
			4	2	2		

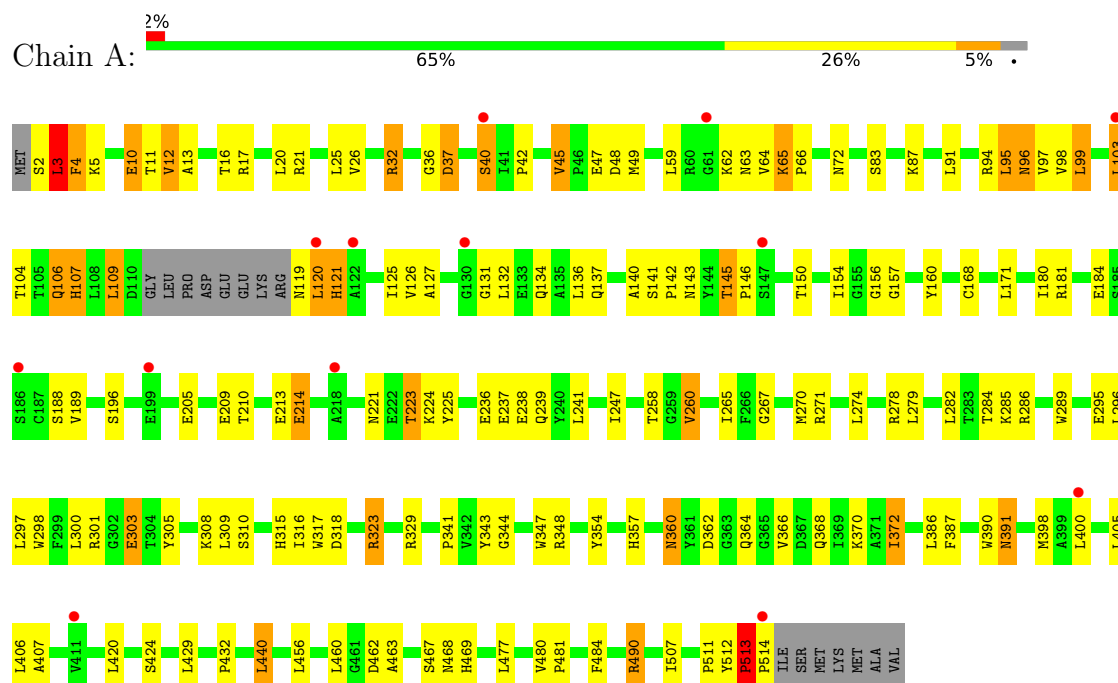
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	20	Total	O	0	0
			20	20		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Bifunctional dihydrofolate reductase-thymidylate synthase (DHFR-TS)



4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	121.05Å 121.05Å 260.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.78 – 3.30 19.78 – 3.30	Depositor EDS
% Data completeness (in resolution range)	97.5 (19.78-3.30) 96.8 (19.78-3.30)	Depositor EDS
R_{merge}	0.01	Depositor
R_{sym}	0.01	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.05 (at 3.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.206 , 0.249 0.199 , 0.241	Depositor DCC
R_{free} test set	738 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	96.7	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 80.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4157	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MTX, SO4, UMP, EDO, NAP, CL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.59	0/4112	0.92	9/5580 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	141	SER	CA-C-N	9.78	132.06	119.84
1	A	141	SER	C-N-CA	9.78	132.06	119.84
1	A	391	ASN	CA-C-N	5.51	125.60	119.32
1	A	391	ASN	C-N-CA	5.51	125.60	119.32
1	A	209	GLU	N-CA-C	-5.44	105.44	112.68
1	A	40	SER	N-CA-C	5.40	118.09	109.02
1	A	45	VAL	CA-C-N	5.02	124.68	119.56
1	A	45	VAL	C-N-CA	5.02	124.68	119.56
1	A	513	PRO	N-CA-C	5.01	116.81	110.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	3	LEU	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4015	0	3966	109	0
2	A	10	0	0	0	0
3	A	3	0	0	0	0
4	A	48	0	25	4	0
5	A	33	0	20	3	0
6	A	20	0	11	1	0
7	A	8	0	12	0	0
8	A	20	0	0	1	0
All	All	4157	0	4034	111	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (111) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:LEU:HA	1:A:5:LYS:H	1.04	1.18
1:A:2:SER:O	1:A:3:LEU:HG	1.48	1.14
1:A:140:ALA:HA	1:A:145:THR:HG23	1.47	0.96
1:A:2:SER:C	1:A:3:LEU:HG	1.93	0.94
1:A:3:LEU:HA	1:A:5:LYS:N	1.81	0.94
1:A:13:ALA:O	1:A:16:THR:HB	1.70	0.92
1:A:120:LEU:HB2	1:A:121:HIS:HB3	1.53	0.90
1:A:32:ARG:HB3	1:A:32:ARG:HH11	1.38	0.87
1:A:91:LEU:HB3	1:A:94:ARG:HH21	1.41	0.82
1:A:119:ASN:O	1:A:120:LEU:HB3	1.84	0.76
1:A:360:ASN:C	1:A:360:ASN:HD22	1.96	0.74
1:A:3:LEU:CA	1:A:5:LYS:H	1.94	0.73
1:A:2:SER:C	1:A:3:LEU:CG	2.62	0.72
1:A:295:GLU:O	1:A:298:TRP:HB3	1.90	0.70
1:A:236:GLU:O	1:A:239:GLN:HG2	1.92	0.69
1:A:214:GLU:OE1	1:A:224:LYS:HD3	1.94	0.68
1:A:120:LEU:CB	1:A:121:HIS:HB3	2.24	0.67
1:A:20:LEU:HD23	1:A:137:GLN:HG3	1.76	0.66
1:A:154:ILE:O	4:A:601:NAP:H5N	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:THR:HG22	1:A:12:VAL:H	1.61	0.65
1:A:221:ASN:HB2	1:A:223:THR:HG22	1.78	0.65
1:A:270:MET:HE2	1:A:460:LEU:HD11	1.77	0.65
1:A:10:GLU:H	1:A:10:GLU:CD	2.06	0.63
1:A:241:LEU:HD11	1:A:284:THR:HG21	1.81	0.63
1:A:103:LEU:HD12	1:A:107:HIS:HD2	1.64	0.62
1:A:131:GLY:H	1:A:134:GLN:NE2	1.99	0.61
1:A:91:LEU:HD21	5:A:602:MTX:H16	1.81	0.60
1:A:131:GLY:N	1:A:134:GLN:HE21	1.99	0.60
1:A:360:ASN:C	1:A:360:ASN:ND2	2.58	0.60
1:A:106:GLN:HA	1:A:109:LEU:HD12	1.85	0.59
1:A:286:ARG:HB3	1:A:514:PRO:HA	1.85	0.58
1:A:140:ALA:HA	1:A:145:THR:CG2	2.30	0.57
1:A:282:LEU:HD12	1:A:432:PRO:HB3	1.87	0.57
1:A:91:LEU:HB3	1:A:94:ARG:NH2	2.16	0.57
1:A:32:ARG:HB3	1:A:32:ARG:NH1	2.15	0.56
1:A:21:ARG:HD2	1:A:507:ILE:HD13	1.86	0.56
1:A:47:GLU:N	1:A:47:GLU:OE1	2.38	0.56
1:A:96:ASN:N	1:A:96:ASN:HD22	2.04	0.56
1:A:11:THR:HG22	1:A:12:VAL:N	2.21	0.55
1:A:121:HIS:O	1:A:121:HIS:ND1	2.40	0.55
1:A:237:GLU:HG2	1:A:481:PRO:HB3	1.88	0.55
1:A:98:VAL:O	1:A:127:ALA:HA	2.06	0.54
1:A:387:PHE:CE1	1:A:407:ALA:HB3	2.42	0.54
1:A:131:GLY:N	1:A:134:GLN:NE2	2.55	0.54
1:A:267:GLY:HA2	1:A:460:LEU:O	2.08	0.54
1:A:3:LEU:N	1:A:4:PHE:HB2	2.23	0.53
1:A:96:ASN:HD22	1:A:96:ASN:H	1.58	0.52
1:A:104:THR:H	1:A:107:HIS:HB2	1.75	0.52
1:A:134:GLN:H	1:A:134:GLN:CD	2.18	0.51
1:A:36:GLY:HA3	1:A:40:SER:O	2.10	0.51
1:A:120:LEU:CA	1:A:121:HIS:HB3	2.41	0.51
1:A:468:ASN:ND2	1:A:469:HIS:CD2	2.78	0.51
4:A:601:NAP:H4N	5:A:602:MTX:NA4	2.27	0.49
1:A:2:SER:HB2	1:A:3:LEU:CD2	2.43	0.49
1:A:45:VAL:HG21	1:A:180:ILE:HD12	1.95	0.49
1:A:48:ASP:OD2	5:A:602:MTX:N1	2.46	0.48
1:A:344:GLY:O	1:A:347:TRP:HB2	2.13	0.48
1:A:120:LEU:O	1:A:120:LEU:HG	2.13	0.48
1:A:59:LEU:HD23	1:A:72:ASN:ND2	2.29	0.48
1:A:258:THR:OG1	1:A:260:VAL:HG23	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:PRO:HG3	1:A:490:ARG:HH21	1.78	0.48
1:A:360:ASN:HD21	1:A:362:ASP:HB2	1.79	0.48
1:A:360:ASN:ND2	1:A:362:ASP:H	2.11	0.47
1:A:21:ARG:H	1:A:136:LEU:HB3	1.79	0.47
1:A:62:LYS:O	1:A:64:VAL:N	2.48	0.47
1:A:424:SER:OG	6:A:603:UMP:H3'	2.15	0.46
1:A:214:GLU:OE1	1:A:224:LYS:CD	2.64	0.46
1:A:512:TYR:HB3	1:A:513:PRO:HD2	1.97	0.46
1:A:47:GLU:HG2	1:A:225:TYR:OH	2.15	0.46
1:A:391:ASN:OD1	1:A:391:ASN:C	2.58	0.46
1:A:301:ARG:HD2	1:A:303:GLU:OE1	2.16	0.46
1:A:196:SER:HB3	1:A:205:GLU:HA	1.98	0.46
1:A:271:ARG:HD2	8:A:906:HOH:O	2.15	0.46
1:A:303:GLU:HG2	1:A:308:LYS:HD3	1.98	0.45
1:A:241:LEU:HD22	1:A:477:LEU:HD23	1.99	0.45
1:A:134:GLN:NE2	1:A:134:GLN:H	2.14	0.45
1:A:62:LYS:O	1:A:64:VAL:HG12	2.17	0.45
1:A:156:GLY:O	1:A:157:GLY:C	2.59	0.45
1:A:316:ILE:HG13	1:A:317:TRP:N	2.32	0.45
1:A:329:ARG:NH1	1:A:398:MET:O	2.50	0.45
1:A:238:GLU:HG3	1:A:481:PRO:HG3	2.00	0.44
1:A:341:PRO:O	1:A:398:MET:HE2	2.18	0.44
1:A:300:LEU:O	1:A:348:ARG:NH1	2.50	0.44
1:A:121:HIS:O	1:A:121:HIS:CG	2.71	0.44
1:A:341:PRO:HG3	1:A:354:TYR:HB2	1.99	0.44
1:A:2:SER:HB2	1:A:3:LEU:HD21	1.98	0.44
1:A:296:LEU:HD22	1:A:440:LEU:HB3	2.00	0.44
1:A:25:LEU:HD11	1:A:160:TYR:CD1	2.53	0.43
1:A:36:GLY:HA2	1:A:42:PRO:HD3	2.00	0.43
1:A:390:TRP:HB2	1:A:405:LEU:HB2	2.00	0.43
4:A:601:NAP:H6N	4:A:601:NAP:H2D	1.57	0.43
1:A:99:LEU:O	4:A:601:NAP:H1B	2.18	0.43
1:A:278:ARG:HH11	1:A:484:PHE:HB2	1.84	0.43
1:A:265:ILE:HG13	1:A:463:ALA:HB3	2.01	0.42
1:A:10:GLU:CD	1:A:10:GLU:N	2.74	0.42
1:A:65:LYS:HD2	1:A:66:PRO:HD2	2.01	0.42
1:A:237:GLU:OE2	1:A:284:THR:HG23	2.19	0.42
1:A:168:CYS:O	1:A:171:LEU:N	2.49	0.42
1:A:305:TYR:CD1	1:A:357:HIS:CE1	3.08	0.42
1:A:323:ARG:HA	1:A:323:ARG:NH1	2.34	0.42
1:A:120:LEU:HA	1:A:121:HIS:HA	1.92	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:468:ASN:HD22	1:A:469:HIS:CD2	2.38	0.42
1:A:462:ASP:OD1	1:A:462:ASP:C	2.63	0.41
1:A:134:GLN:CD	1:A:134:GLN:N	2.78	0.41
1:A:368:GLN:O	1:A:372:ILE:HG12	2.20	0.41
1:A:247:ILE:HG12	1:A:265:ILE:HG23	2.02	0.41
1:A:278:ARG:NH1	1:A:484:PHE:HD2	2.17	0.41
1:A:279:LEU:HD21	1:A:289:TRP:CZ3	2.56	0.41
1:A:37:ASP:HA	1:A:188:SER:H	1.86	0.40
1:A:315:HIS:HA	1:A:318:ASP:OD1	2.21	0.40
1:A:95:LEU:HD22	1:A:97:VAL:HG23	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	501/521 (96%)	445 (89%)	45 (9%)	11 (2%)	5	26

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	PHE
1	A	120	LEU
1	A	142	PRO
1	A	343	TYR
1	A	37	ASP
1	A	63	ASN
1	A	121	HIS
1	A	513	PRO
1	A	109	LEU
1	A	3	LEU
1	A	511	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	431/446 (97%)	379 (88%)	52 (12%)	5 20

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	10	GLU
1	A	12	VAL
1	A	17	ARG
1	A	26	VAL
1	A	32	ARG
1	A	49	MET
1	A	65	LYS
1	A	83	SER
1	A	87	LYS
1	A	95	LEU
1	A	96	ASN
1	A	99	LEU
1	A	103	LEU
1	A	106	GLN
1	A	107	HIS
1	A	125	ILE
1	A	126	VAL
1	A	132	LEU
1	A	143	ASN
1	A	145	THR
1	A	150	THR
1	A	181	ARG
1	A	184	GLU
1	A	189	VAL
1	A	210	THR
1	A	213	GLU
1	A	214	GLU
1	A	223	THR
1	A	260	VAL

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Mol	Chain	Res	Type
1	A	274	LEU
1	A	285	LYS
1	A	297	LEU
1	A	303	GLU
1	A	309	LEU
1	A	310	SER
1	A	323	ARG
1	A	360	ASN
1	A	364	GLN
1	A	366	VAL
1	A	370	LYS
1	A	372	ILE
1	A	386	LEU
1	A	400	LEU
1	A	406	LEU
1	A	420	LEU
1	A	429	LEU
1	A	440	LEU
1	A	456	LEU
1	A	467	SER
1	A	480	VAL
1	A	490	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	HIS
1	A	134	GLN
1	A	137	GLN
1	A	320	ASN
1	A	346	GLN
1	A	360	ASN
1	A	364	GLN
1	A	422	GLN
1	A	468	ASN
1	A	469	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 3 are monoatomic - leaving 7 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAP	A	601	-	50,52,52	1.78	7 (14%)	71,80,80	1.64	11 (15%)
7	EDO	A	701	-	3,3,3	0.42	0	2,2,2	0.55	0
7	EDO	A	702	-	3,3,3	0.46	0	2,2,2	0.36	0
6	UMP	A	603	-	21,21,21	1.08	2 (9%)	30,31,31	1.88	8 (26%)
2	SO4	A	801	-	4,4,4	0.29	0	6,6,6	0.17	0
5	MTX	A	602	-	35,35,35	1.88	4 (11%)	47,49,49	1.78	9 (19%)
2	SO4	A	802	-	4,4,4	0.27	0	6,6,6	0.23	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAP	A	601	-	-	12/35/67/67	0/5/5/5
7	EDO	A	701	-	-	1/1/1/1	-
7	EDO	A	702	-	-	1/1/1/1	-
6	UMP	A	603	-	-	6/10/22/22	0/2/2/2
5	MTX	A	602	-	-	11/25/25/25	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	601	NAP	O7N-C7N	9.25	1.41	1.24
5	A	602	MTX	O-C	8.04	1.42	1.23
5	A	602	MTX	C7-N8	4.26	1.39	1.31
5	A	602	MTX	C6-N5	3.81	1.39	1.32
4	A	601	NAP	C2N-N1N	3.43	1.38	1.35
4	A	601	NAP	C2A-N3A	2.79	1.38	1.33
4	A	601	NAP	C2A-N1A	2.72	1.38	1.33
6	A	603	UMP	C2-N1	2.40	1.42	1.38
5	A	602	MTX	C4-N3	2.35	1.38	1.33
4	A	601	NAP	C8A-N7A	2.14	1.35	1.31
6	A	603	UMP	C6-C5	2.08	1.39	1.35
4	A	601	NAP	PA-O3	2.02	1.61	1.59
4	A	601	NAP	P2B-O2B	2.01	1.63	1.59

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	601	NAP	N3A-C2A-N1A	-5.74	119.89	128.58
5	A	602	MTX	N1-C2-N3	-5.42	120.33	127.21
4	A	601	NAP	C5A-C4A-N3A	-5.04	119.78	126.72
6	A	603	UMP	C4-N3-C2	-4.39	121.16	126.61
4	A	601	NAP	C2B-C1B-N9A	-4.39	106.53	113.75
5	A	602	MTX	N8-C8A-N1	4.13	120.27	115.77
5	A	602	MTX	C2-N1-C8A	4.05	119.85	115.48
6	A	603	UMP	C1'-N1-C2	4.04	125.55	117.66
4	A	601	NAP	C2A-N3A-C4A	3.70	120.87	111.83
6	A	603	UMP	C5-C4-N3	3.64	119.89	114.80
4	A	601	NAP	N9A-C8A-N7A	-3.43	109.07	113.94
4	A	601	NAP	N3A-C4A-N9A	3.42	132.98	127.17
4	A	601	NAP	C5A-N7A-C8A	3.36	108.73	103.45
6	A	603	UMP	C1'-N1-C6	-3.20	115.24	121.53
4	A	601	NAP	C6N-N1N-C2N	-3.13	119.21	121.88
5	A	602	MTX	C6-C7-N8	-3.10	120.16	123.14
6	A	603	UMP	N3-C2-N1	3.10	118.92	114.89
6	A	603	UMP	O4-C4-C5	-3.08	119.85	125.16
6	A	603	UMP	OP3-P-O5'	2.96	114.39	106.67
5	A	602	MTX	C7-C6-N5	-2.75	119.09	120.87
5	A	602	MTX	CB-CG-CD	-2.49	105.85	112.49
6	A	603	UMP	O2-C2-N3	-2.45	116.97	121.49
5	A	602	MTX	C8A-C4A-N5	-2.44	119.62	122.35
4	A	601	NAP	C3N-C7N-N7N	2.35	120.63	117.74
4	A	601	NAP	C4A-C5A-N7A	-2.31	107.94	110.58
4	A	601	NAP	C4D-O4D-C1D	2.12	111.86	109.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	602	MTX	C7-N8-C8A	2.11	120.14	117.20
5	A	602	MTX	C6-N5-C4A	2.01	121.01	118.17

There are no chirality outliers.

All (31) torsion outliers are listed below:

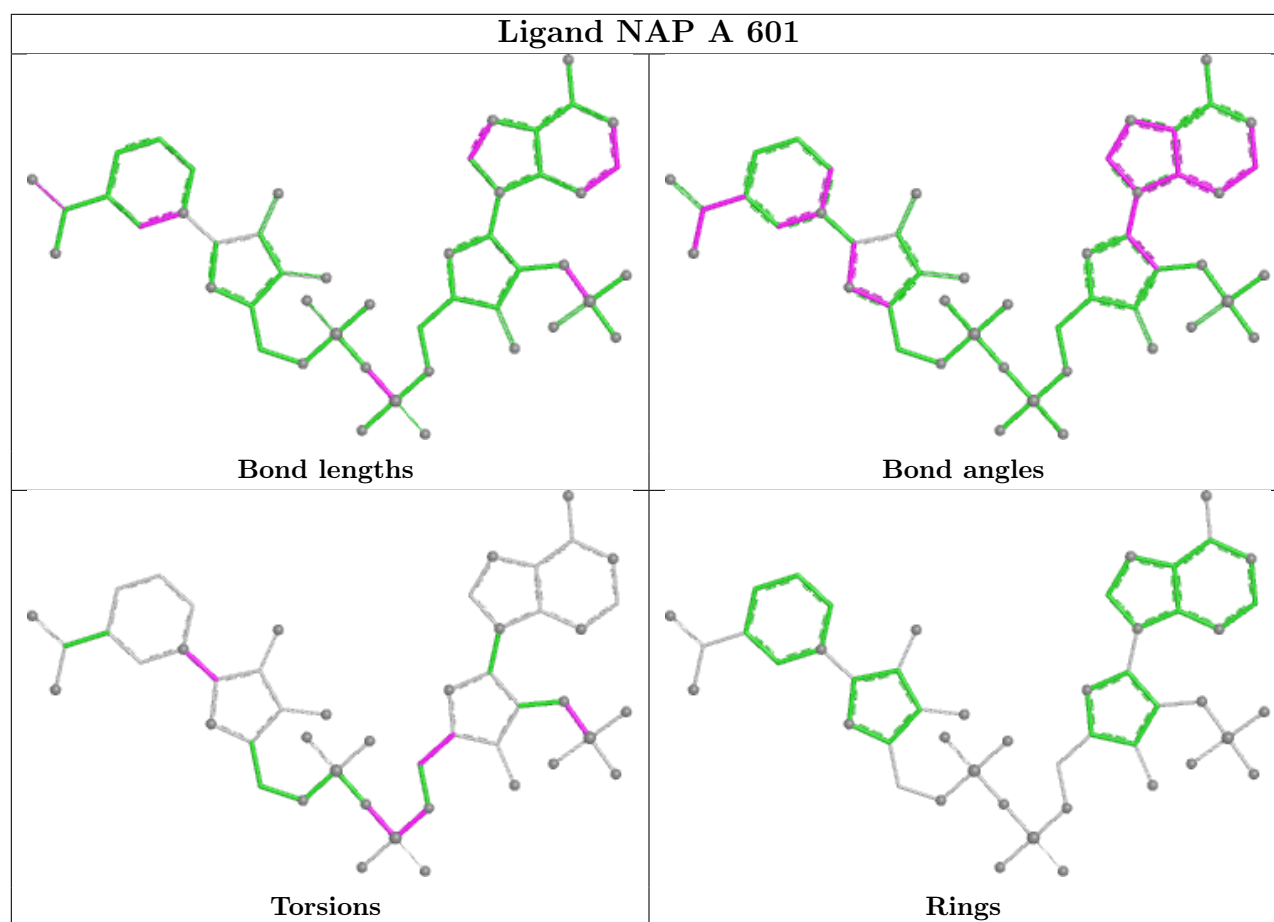
Mol	Chain	Res	Type	Atoms
4	A	601	NAP	C5B-O5B-PA-O1A
4	A	601	NAP	C5B-O5B-PA-O2A
4	A	601	NAP	C5B-O5B-PA-O3
4	A	601	NAP	O4B-C4B-C5B-O5B
4	A	601	NAP	O4D-C1D-N1N-C2N
4	A	601	NAP	O4D-C1D-N1N-C6N
4	A	601	NAP	C2D-C1D-N1N-C2N
4	A	601	NAP	C2D-C1D-N1N-C6N
5	A	602	MTX	N-CA-CB-CG
5	A	602	MTX	CB-CA-N-C
6	A	603	UMP	C2'-C1'-N1-C2
4	A	601	NAP	C3B-C4B-C5B-O5B
6	A	603	UMP	C3'-C4'-C5'-O5'
6	A	603	UMP	O4'-C4'-C5'-O5'
6	A	603	UMP	C2'-C1'-N1-C6
5	A	602	MTX	CT-CA-CB-CG
7	A	701	EDO	O1-C1-C2-O2
5	A	602	MTX	C11-C-N-CA
5	A	602	MTX	O-C-N-CA
5	A	602	MTX	C13-C14-N10-CM
4	A	601	NAP	PN-O3-PA-O5B
5	A	602	MTX	C6-C9-N10-CM
5	A	602	MTX	C15-C14-N10-CM
5	A	602	MTX	C13-C14-N10-C9
4	A	601	NAP	C2B-O2B-P2B-O1X
6	A	603	UMP	C4'-C5'-O5'-P
5	A	602	MTX	OE2-CD-CG-CB
7	A	702	EDO	O1-C1-C2-O2
6	A	603	UMP	C5'-O5'-P-OP2
5	A	602	MTX	OE1-CD-CG-CB
4	A	601	NAP	C2B-O2B-P2B-O2X

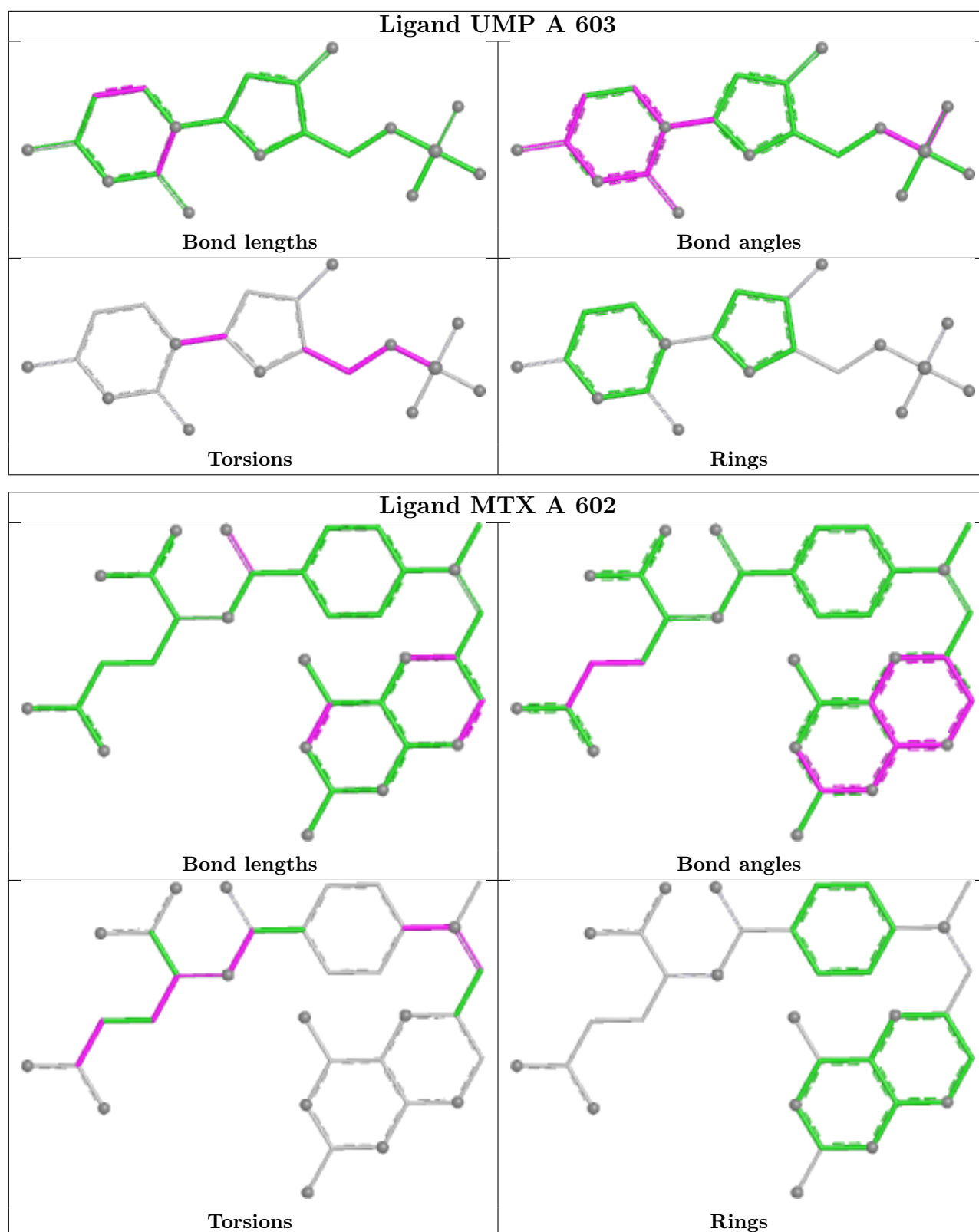
There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	601	NAP	4	0
6	A	603	UMP	1	0
5	A	602	MTX	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	505/521 (96%)	0.28	13 (2%) 57 38	73, 89, 103, 114	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	61	GLY	2.7
1	A	400	LEU	2.5
1	A	186	SER	2.4
1	A	122	ALA	2.3
1	A	147	SER	2.3
1	A	411	VAL	2.2
1	A	103	LEU	2.2
1	A	130	GLY	2.2
1	A	514	PRO	2.2
1	A	218	ALA	2.2
1	A	40	SER	2.1
1	A	120	LEU	2.1
1	A	199	GLU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

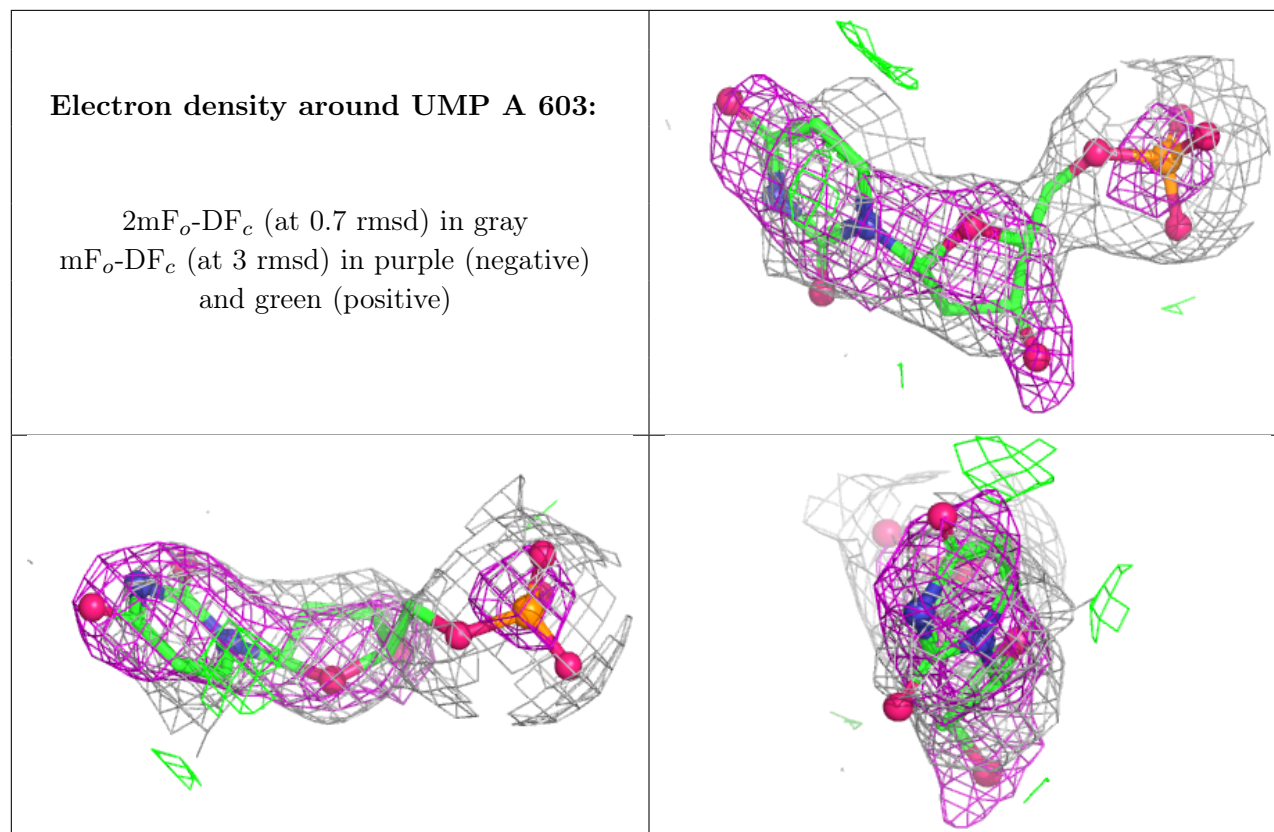
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

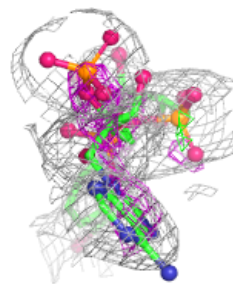
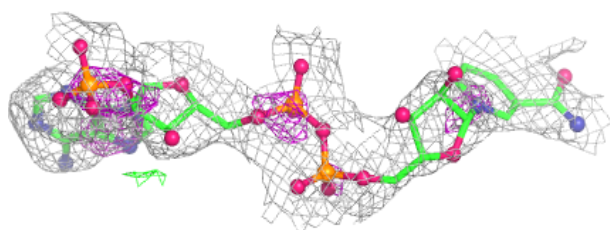
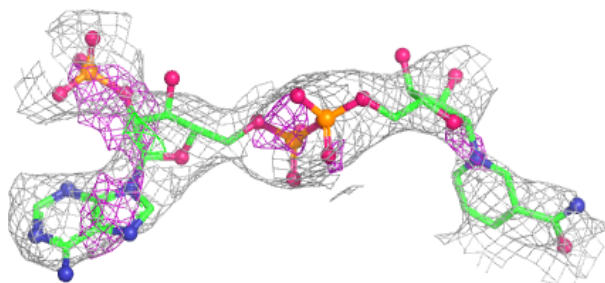
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	UMP	A	603	20/20	0.78	0.14	80,101,106,107	0
2	SO4	A	801	5/5	0.82	0.13	124,124,124,125	0
4	NAP	A	601	48/48	0.83	0.14	123,126,130,131	0
7	EDO	A	702	4/4	0.84	0.10	87,87,87,88	0
3	CL	A	903	1/1	0.86	0.10	98,98,98,98	0
3	CL	A	902	1/1	0.87	0.08	117,117,117,117	0
7	EDO	A	701	4/4	0.88	0.16	79,80,80,80	0
5	MTX	A	602	33/33	0.90	0.12	98,98,99,99	0
2	SO4	A	802	5/5	0.94	0.22	109,110,110,111	0
3	CL	A	901	1/1	0.97	0.07	99,99,99,99	0

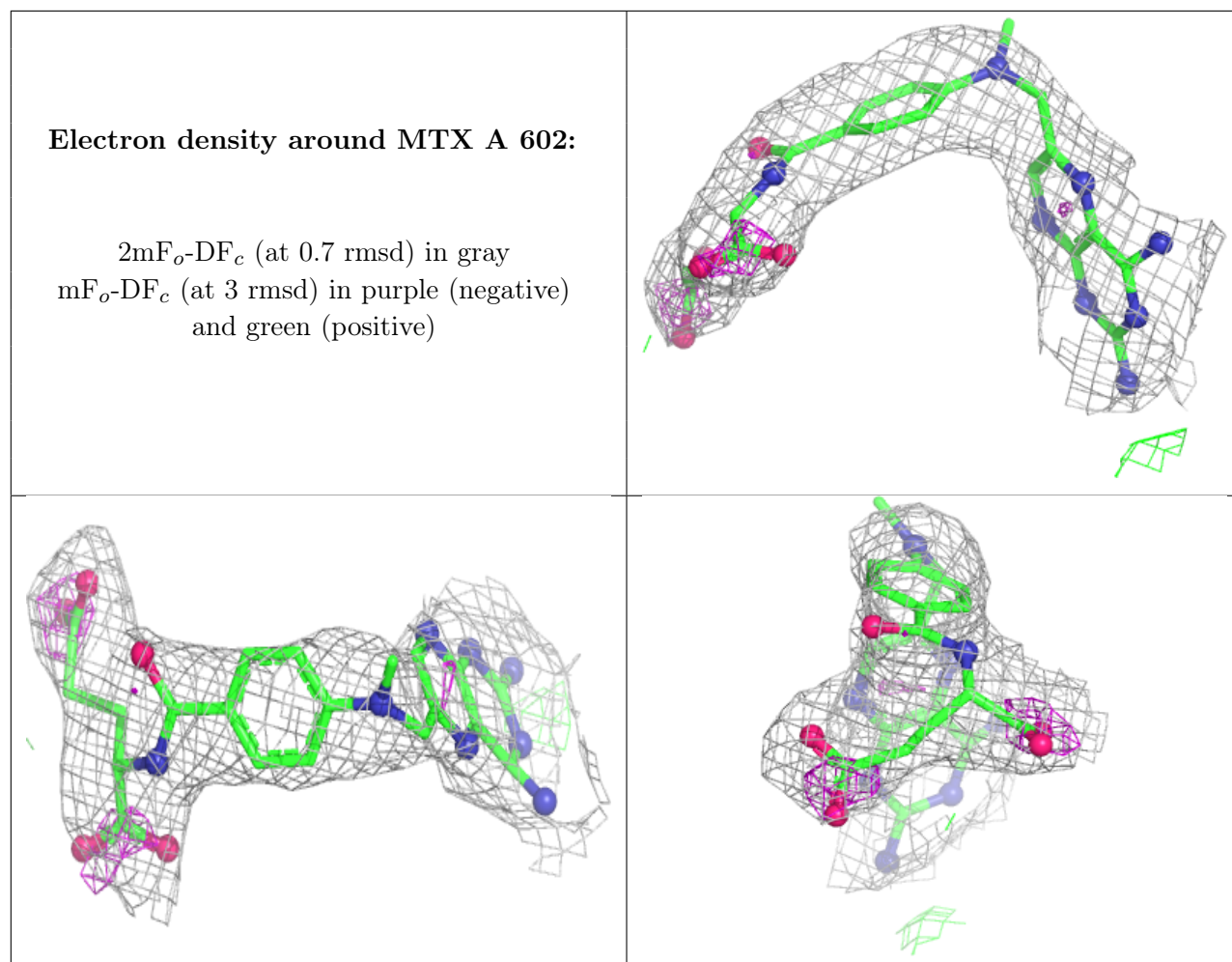
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around NAP A 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers ⓘ

There are no such residues in this entry.